

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
 Data File : BM048129.D
 Acq On : 18 Oct 2024 10:46
 Operator : RC/JU
 Sample : P4363-01
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EOAY5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048129.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	38	42	50	rVB	61999	92136	1.84%	0.182%
2	3.969	165	167	174	rVB2	15134	19955	0.40%	0.040%
3	4.134	192	195	201	rVB7	7448	12349	0.25%	0.024%
4	4.346	227	231	238	rVB5	13435	22416	0.45%	0.044%
5	4.469	249	252	258	rBV4	7809	11505	0.23%	0.023%
6	4.528	258	262	267	rVB	28389	40327	0.81%	0.080%
7	4.769	299	303	308	rVB7	8637	12386	0.25%	0.025%
8	4.893	320	324	330	rVB4	10471	18762	0.37%	0.037%
9	5.581	435	441	446	rBV8	8800	18620	0.37%	0.037%
10	5.781	469	475	477	rBV	24360	37515	0.75%	0.074%
11	6.034	513	518	524	rVB5	11345	18977	0.38%	0.038%
12	6.151	529	538	547	rVB2	82639	167378	3.34%	0.332%
13	6.581	606	611	614	rBV7	9445	13519	0.27%	0.027%
14	6.628	614	619	625	rVB3	17558	33471	0.67%	0.066%
15	6.793	643	647	651	rBV6	4565	7688	0.15%	0.015%
16	6.946	662	673	686	rBV	420971	765403	15.28%	1.516%
17	7.098	693	699	713	rBV	780992	1304460	26.04%	2.584%
18	7.304	725	734	747	rBV	1009727	1686176	33.67%	3.340%
19	7.769	806	813	822	rBV	789526	1319244	26.34%	2.613%
20	8.481	927	934	952	rBV	963724	1638650	32.72%	3.245%
21	8.922	1002	1009	1021	rBV	866390	1502846	30.01%	2.976%
22	9.351	1075	1082	1087	rBV	17205	30027	0.60%	0.059%
23	9.645	1122	1132	1139	rBV	716883	1239882	24.76%	2.456%
24	9.975	1186	1188	1195	rVB4	4335	6855	0.14%	0.014%
25	10.186	1217	1224	1244	rBV	918873	1894879	37.83%	3.753%
26	10.557	1280	1287	1302	rBV	1054168	1801870	35.98%	3.569%
27	10.710	1305	1313	1333	rVB2	86098	244241	4.88%	0.484%
28	11.839	1498	1505	1509	rBV2	17043	28887	0.58%	0.057%
29	11.992	1524	1531	1535	rVB8	2769	6311	0.13%	0.012%
30	12.175	1556	1562	1570	rBV3	11880	25286	0.50%	0.050%
31	12.404	1594	1601	1604	rBV3	6772	9620	0.19%	0.019%
32	12.769	1659	1663	1666	rBV6	3575	5514	0.11%	0.011%
33	13.022	1701	1706	1711	rVB6	5314	10078	0.20%	0.020%
34	13.816	1834	1841	1857	rBV	1792089	2618056	52.27%	5.185%
35	14.098	1882	1889	1903	rBV	2182715	3245897	64.81%	6.429%
36	14.404	1934	1941	1953	rBV	1484672	2231068	44.55%	4.419%

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Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
 Title : SVOA CALIBRATION

37	14.627	1973	1979	2000	rBV	167071	474960	9.48%	0.941%
38	15.204	2071	2077	2079	rBV2	10787	13820	0.28%	0.027%
39	15.398	2103	2110	2125	rBV	2623673	3951694	78.90%	7.827%
40	15.521	2125	2131	2146	rVV	822536	1251753	24.99%	2.479%
41	15.639	2148	2151	2159	rVB2	14352	24583	0.49%	0.049%
42	15.763	2169	2172	2180	rVB8	4342	9178	0.18%	0.018%
43	15.968	2204	2207	2211	rVB5	4426	5846	0.12%	0.012%
44	16.104	2223	2230	2241	rBV	45334	120203	2.40%	0.238%
45	16.174	2241	2242	2246	rVB3	6074	7145	0.14%	0.014%
46	16.833	2349	2354	2356	rVB5	3919	6194	0.12%	0.012%
47	16.933	2366	2371	2374	rVB5	4151	5234	0.10%	0.010%
48	17.145	2399	2407	2417	rBV	1710462	2482324	49.56%	4.916%
49	17.245	2417	2424	2442	rVV	2800566	4260661	85.07%	8.438%
50	17.433	2454	2456	2460	rVB4	4317	5132	0.10%	0.010%
51	17.545	2471	2475	2477	rBV4	3688	5596	0.11%	0.011%
52	17.810	2514	2520	2526	rBV9	8654	17464	0.35%	0.035%
53	17.980	2544	2549	2551	rBV5	3587	5134	0.10%	0.010%
54	18.039	2555	2559	2568	rVV8	23010	48981	0.98%	0.097%
55	18.115	2568	2572	2576	rVB3	9788	15315	0.31%	0.030%
56	18.251	2590	2595	2596	rBV5	4146	6398	0.13%	0.013%
57	18.939	2710	2712	2717	rBV5	11672	12995	0.26%	0.026%
58	19.098	2736	2739	2746	rBV2	35822	46976	0.94%	0.093%
59	19.168	2748	2751	2755	rBV	32537	44956	0.90%	0.089%
60	19.315	2774	2776	2783	rBV7	13083	20931	0.42%	0.041%
61	19.533	2807	2813	2829	rBV	3605246	4990188	99.63%	9.883%
62	21.368	3120	3125	3139	rBV	1593332	2618112	52.27%	5.185%
63	24.144	3589	3597	3616	rVB	1946378	5008531	100.00%	9.920%
64	24.350	3624	3632	3647	rVB	1096067	2888404	57.67%	5.721%

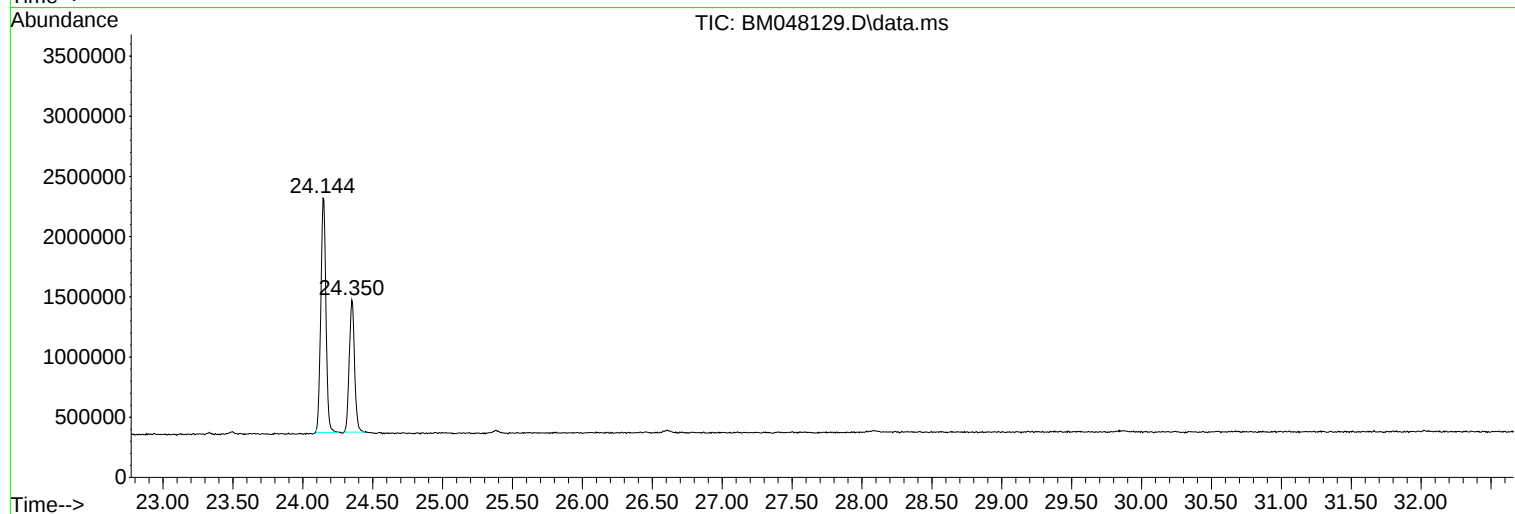
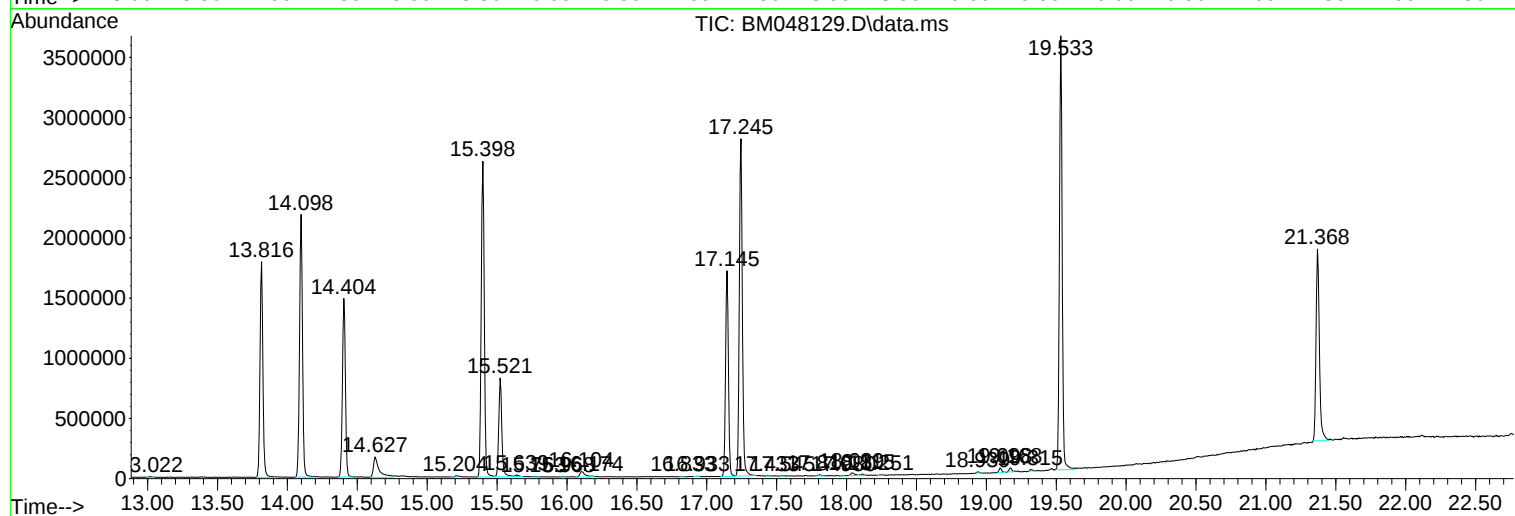
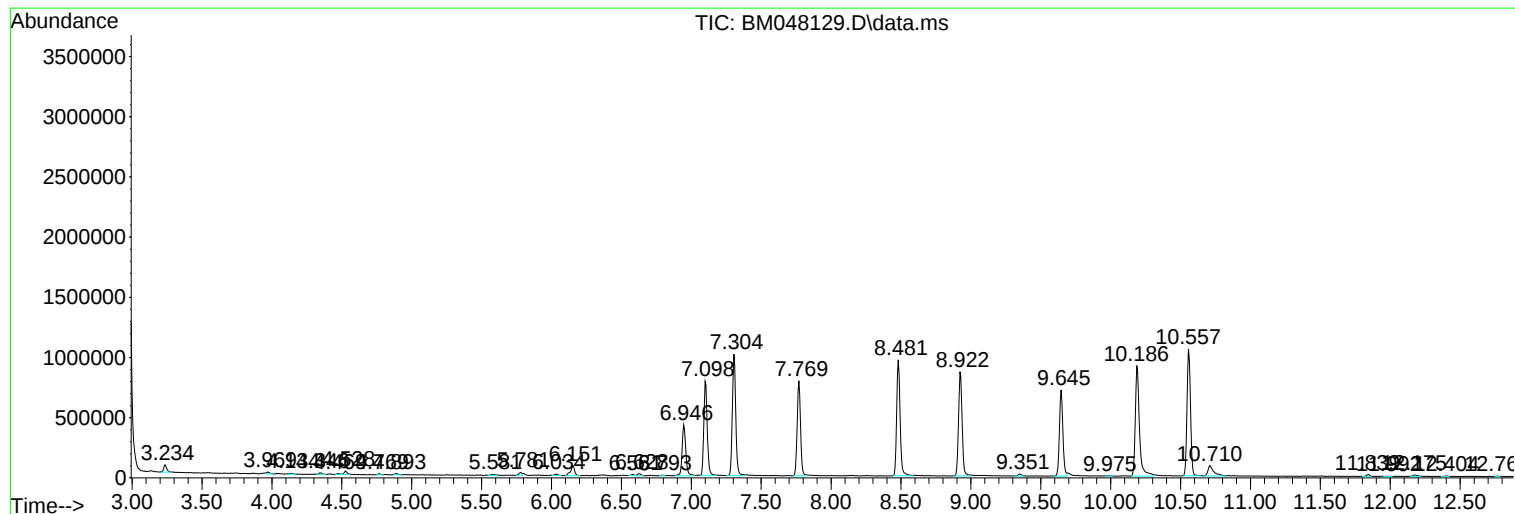
Sum of corrected areas: 50490962

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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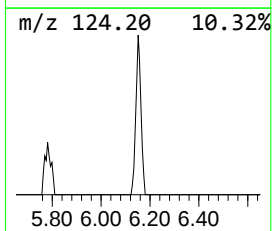
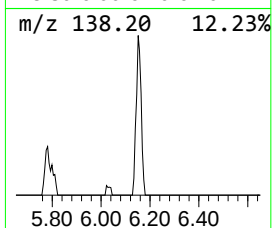
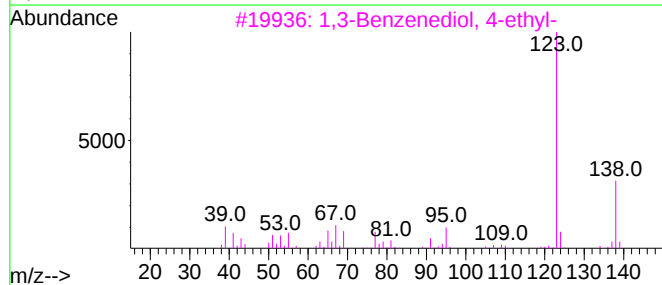
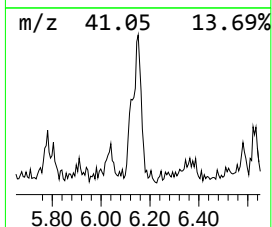
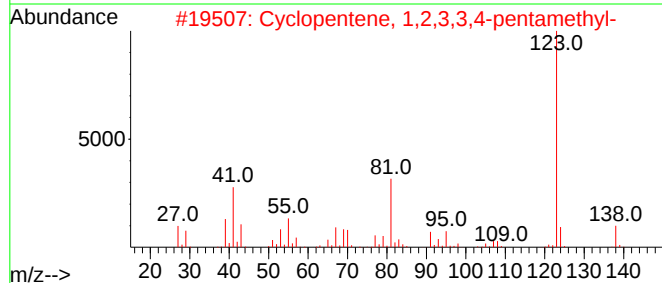
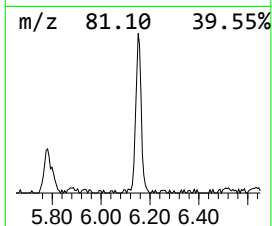
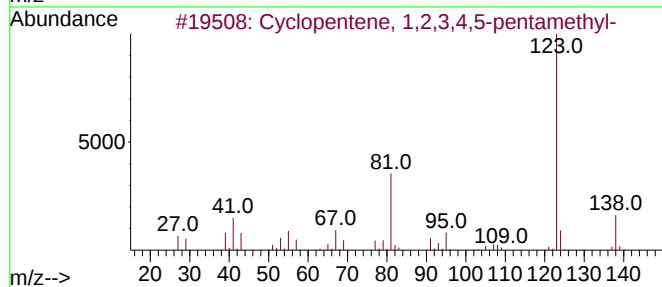
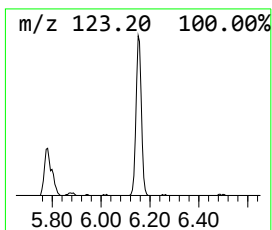
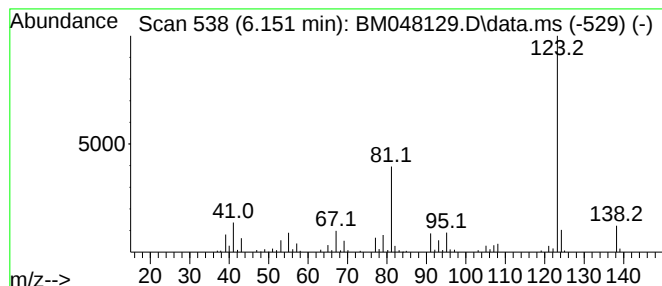
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclopentene, 1,2,3,4,5-pen... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.151	2.54 ng/ul	167378	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentene, 1,2,3,4,5-pentamet...	138	C10H18	1000154-28-6	91
2			Cyclopentene, 1,2,3,3,4-pentamet...	138	C10H18	197390-29-7	87
3			1,3-Benzenediol, 4-ethyl-	138	C8H10O2	002896-60-8	72
4			1,3-Benzenediol, 4-ethyl-	138	C8H10O2	002896-60-8	72
5			1,3-Benzenediol, 4-ethyl-	138	C8H10O2	002896-60-8	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopentene, 1...	6.151	2.5	ng/ul	167378	1	7.769	1319240	20.0